In the Claims:

1-15. (Canceled)

16. (Previously Presented) A compound of formula (I)

wherein

---- represents a single or a double bond;

R is a radical selected from:

in which R₁ is halogen, cyano, C₁₋₄ alkyl, C₁₋₄ alkoxy, trifluoromethyl or trifluoromethoxy and p is zero or an integer from 1 to 3;

R₂ is hydrogen or C₁₋₄ alkyl;

R₃ is hydrogen, hydroxy or C₁₋₄ alkyl;

 R_4 is hydrogen or R_4 together with R_3 is =0 or = CH_2 ;

R₅ is phenyl, naphthyl, a 9 to 10 membered fused bicyclic heterocyclic group or a 5 or 6 membered heteroaryl group, wherein said groups are optionally substituted by 1 to 3 groups independently selected from trifluoromethyl, C₁₋₄ alkyl, hydroxy, cyano, C₁₋₄ alkoxy, trifluoromethoxy, halogen or S(O)_qC₁₋₄ alkyl;

R₆ and R₇ independently are hydrogen, cyano, C₁₋₄ alkyl;

R₈ is (CH₂)_rR₁₀;

Rg is hydrogen, halogen, C₃₋₇ cycloalkyl, hydroxy, nitro, cyano or C₁₋₄ alkyl optionally substituted by one or two groups selected from halogen, cyano, hydroxy or C₁₋₄ alkoxy;

R₁₀ is hydrogen or C₃₋₇ cycloalkyl;

n is 2;

q is 0, 1 or 2;

r is 0 or an integer from 1 to 4;

or a pharmaceutically acceptable salt thereof.

- 17. (Canceled).
- 18. (Previously Presented) A compound as claimed in claim 1 wherein R is:

wherein R₁ is halogen, C₁₋₄ alkyl, cyano, C₁₋₄ alkoxy, trifluoromethyl or trifluoromethoxy and p is zero or an integer from 1 to 3.

- 19. (Previously Presented) A compound as claimed in claim 1 wherein R_5 is phenyl or naphthyl optionally substituted by one or two groups selected from trifluoromethyl, cyano, C_{1-4} alkyl or halogen.
- 20. (Previously Presented) A compound as claimed claim 1 wherein R_8 is $(CH_2)_rR_{10}$ wherein R_{10} is hydrogen or C_{3-7} cycloalkyl and r is 0 or 1.
- 21. (Previously Presented) A compound as claimed in claim 1, wherein R₉ is hydrogen or C₁₋₄ alkyl optionally substituted by one or two halogens.
- 22. (Previously Presented) A compound as claimed in claim 1 wherein:

R is phenyl substituted by a fluorine;

R₂, R₉ and R₄ are each hydrogen;

 R_3 is hydrogen, hydroxy or methyl, or R_3 together with R_4 is =0 or = CH_2 ;

R6 and R7 are independently hydrogen or methyl;

 $\ensuremath{\mathsf{R}}_5$ is phenyl or naphthyl optionally substituted by one or two groups independently selected from $% \ensuremath{\mathsf{cyano}}$ cyano, methyl, chlorine, bromine or fluorine atom; and

Rg is hydrogen, methyl or cyclopropylmethyl.

- 23. (Previously Presented) A compound selected from:
- 1-[(3,5-Dichlorophenyl)methyl]-3-[4-(4-fluorophenyl)-4-piperidinyl]-1,5-dihydro-2*H*-pyrrol-2-one;
- 1-[(3,5-Dichlorophenyl)methyl]-3-[4-(4-fluorophenyl)-1-methyl-4-piperidinyl]-1,5-dihydro-2*H*-pyrrol-2-one;
- 1-[1-(3,5-Dichlorophenyl)ethyl]-3-[4-(4-fluorophenyl)-4-piperidinyl]-1,5-dihydro-2*H*-pyrrol-2-one (Chain Enantiomer 1);
- 1-[(1*S*)-1-(3-Chloro-1-naphthalenyl)ethyl]-3-[4-(4-fluorophenyl)-4-piperidinyl]-1,5-dihydro-2*H*-pyrrol-2-one;
- 1-[(3-Chloro-1-naphthalenyl)methyl]-3-[4-(4-fluorophenyl)-4-piperidinyl]-1,5-dihydro-2*H*-pyrrol-2-one;
- 4-({3-[4-(4-Fluorophenyl)-4-piperidinyl]-2-oxo-2,5-dihydro-1*H*-pyrrol-1-yl}methyl)-2-naphthalenecarbonitrile;
- 1-[1-(3,5-Dichlorophenyl)ethyl]-3-[4-(4-fluorophenyl)-4-piperidinyl]-1,5-dihydro-2*H*-pyrrol-2-one (Chain Enantiomer 2);
- 1-[(1*R*)-1-(3-Chloro-1-naphthalenyl)ethyl]-3-[4-(4-fluorophenyl)-4-piperidinyl]-1,5-dihydro-2*H*-pyrrol-2-one;
- 1-[1-(3,5-Dichlorophenyl)ethyl]-3-[4-(4-fluorophenyl)-1-methyl-4-piperidinyl]-1,5-dihydro-2*H*-pyrrol-2-one (Chain Enantiomer 1);
- 1-[(1*S*)-1-(3-Chloro-1-naphthalenyl)ethyl]-3-[4-(4-fluorophenyl)-1-methyl-4-piperidinyl]-1,5-dihydro-2*H*-pyrrol-2-one;

- 1-[(3-Chloro-1-naphthalenyl)methyl]-3-[4-(4-fluorophenyl)-1-methyl-4-piperidinyl]-1,5-dihydro-2*H*-pyrrol-2-one;
- 1-[1-(3,5-Dichlorophenyl)ethyl]-3-[4-(4-fluorophenyl)-1-methyl-4-piperidinyl]-1,5-dihydro-2*H*-pyrrol-2-one (Chain Enantiomer 2);
- 1-[(1*R*)-1-(3-Chloro-1-naphthalenyl)ethyl]-3-[4-(4-fluorophenyl)-1-methyl-4-piperidinyl]-1,5-dihydro-2*H*-pyrrol-2-one;
- 4-({3-[4-(4-Fluorophenyl)-1-methyl-4-piperidinyl]-2-oxo-2,5-dihydro-1*H*-pyrrol-1-vl}methyl)-2-naphthalenecarbonitrile;
- 1-[1-(3,5-Dichlorophenyl)ethyl]-3-[4-(4-fluorophenyl)-1-methyl-4-piperidinyl]-1,5-dihydro-2*H*-pyrrol-2-one hydrochloride (Chain Enantiomer 1);
- 1-[1-(3,5-Dichlorophenyl)ethyl]-3-[4-(4-fluorophenyl)-1-methyl-4-piperidinyl]-1,5-dihydro-2*H*-pyrrol-2-one hydrochloride (Chain Enantiomer 1);
- 1-[(3-Chloro-1-naphthalenyl)methyl]-3-[1-(cyclopropylmethyl)-4-(4-fluorophenyl)-4-piperidinyl]-1,5-dihydro-2*H*-pyrrol-2-one;
- 1-[(3,5-Dichlorophenyl)methyl]-3-[4-(4-fluorophenyl)-4-piperidinyl]-2-pyrrolidinone
- 1-[(1*S*)-1-(3-Chloro-1-naphthalenyl)ethyl]-3-[4-(4-fluorophenyl)-4-piperidinyl]-2-pyrrolidinone (Diastereoisomer 1);
- 1-[(1*S*)-1-(3-Chloro-1-naphthalenyl)ethyl]-3-[4-(4-fluorophenyl)-4-piperidinyl]-2-pyrrolidinone;
- 1-[1-(3,5-Dichlorophenyl)ethyl]-3-[4-(4-fluorophenyl)-4-piperidinyl]-2-pyrrolidinone (Diastereoisomer 1 Chain Enantiomer 1);
- 1-[1-(3,5-Dichlorophenyl)ethyl]-3-[4-(4-fluorophenyl)-4-piperidinyl]-2-pyrrolidinone (Diastereoisomer 2 Chain Enantiomer 1);
- 1-[1-(3,5-Dichlorophenyl)ethyl]-3-[4-(4-fluorophenyl)-4-piperidinyl]-2-pyrrolidinone (Diastereoisomer 1 Chain Enantiomer 2);
- 1-[1-(3,5-Dichlorophenyl)ethyl]-3-[4-(4-fluorophenyl)-4-piperidinyl]-2-pyrrolidinone (Diastereoisomer 2 Chain Enantiomer 2);
- 4-({3-[4-(4-Fluorophenyl)-4-piperidinyl]-2-oxo-1-pyrrolidinyl}methyl)-2-naphthalenecarbonitrile (Enantiomer 1);
- 4-({3-[4-(4-Fluorophenyl)-4-piperidinyl]-2-oxo-1-pyrrolidinyl}methyl)-2-naphthalenecarbonitrile (Enantiomer 2);

- 7-Fluoro-4-({3-[4-(4-fluorophenyl)-4-piperidinyl]-2-oxo-1-pyrrolidinyl}methyl)-2-naphthalenecarbonitrile (Enantiomer 2);
- 6-Fluoro-4-({3-[4-(4-fluorophenyl)-4-piperidinyl]-2-oxo-1-pyrrolidinyl}methyl)-2-naphthalenecarbonitrile (Enantiomer 2);
- 7-Fluoro-4-({3-[4-(4-fluorophenyl)-4-piperidinyl]-2-oxo-1-pyrrolidinyl}methyl)-2-naphthalenecarbonitrile (Enantiomer 1);
- 6-Fluoro-4-({3-[4-(4-fluorophenyl)-4-piperidinyl]-2-oxo-1-pyrrolidinyl}methyl)-2-naphthalenecarbonitrile (Enantiomer 1);
- 1-[(3,5-Dichlorophenyl)methyl]-3-[4-(4-fluorophenyl)-1-methyl-4-piperidinyl]-2-pyrrolidinone;
- 1-[1-(3-Chloro-1-naphthalenyl)ethyl]-3-[4-(4-fluorophenyl)-1-methyl-4-piperidinyl]2-pyrrolidinone (Diastereoisomer 2 Chain Enantiomer 1);
- 1-[1-(3,5-Dichlorophenyl)ethyl]-3-[4-(4-fluorophenyl)-1-methyl-4-piperidinyl]-2-pyrrolidinone (Diastereoisomer 1 Chain Enantiomer 1);
- 1-[1-(3,5-Dichlorophenyl)ethyl]-3-[4-(4-fluorophenyl)-1-methyl-4-piperidinyl]-2-pyrrolidinone (Diastereoisomer 2 Chain Enantiomer 1);
- 1-[1-(3,5-Dichlorophenyl)ethyl]-3-[4-(4-fluorophenyl)-1-methyl-4-piperidinyl]-2-pyrrolidinone (Diastereoisomer 1 Chain Enantiomer 2);
- 1-[1-(3,5-Dichlorophenyl)ethyl]-3-[4-(4-fluorophenyl)-1-methyl-4-piperidinyl]-2-pyrrolidinone (Diastereoisomer 2 Chain Enantiomer 2);
- 1-[(3,5-Dichlorophenyl)methyl]-3-[4-(4-fluorophenyl)-1-methyl-4-piperidinyl]-2-pyrrolidinone (Enantiomer 1);
- 1-[(3,5-Dichlorophenyl)methyl]-3-[4-(4-fluorophenyl)-1-methyl-4-piperidinyl]-2-pyrrolidinone (Enantiomer 2);
- 4-({3-[4-(4-Fluorophenyl)-1-methyl-4-piperidinyl]-2-oxo-1-pyrrolidinyl}methyl)-2-naphthalenecarbonitrile (Enantiomer 1);
- 4-({3-[4-(4-Fluorophenyl)-1-methyl-4-piperidinyl]-2-oxo-1-pyrrolidinyl}methyl)-2-naphthalenecarbonitrile (Enantiomer 2);
- 7-Fluoro-4-({3-[4-(4-fluorophenyl)-1-methyl-4-piperidinyl]-2-oxo-1-pyrrolidinyl}methyl)-2-naphthalenecarbonitrile (Enantiomer 2);
- 6-Fluoro-4-({3-[4-(4-fluorophenyl)-1-methyl-4-piperidinyl]-2-oxo-1-pyrrolidinyl}methyl)-2-naphthalenecarbonitrile;

- 7-Fluoro-4-({3-[4-(4-fluorophenyl)-1-methyl-4-piperidinyl]-2-oxo-1-pyrrolidinyl}methyl)-2-naphthalenecarbonitrile (Enantiomer 1);
- 6-Fluoro-4-({3-[4-(4-fluorophenyl)-1-methyl-4-piperidinyl]-2-oxo-1-pyrrolidinyl}methyl)-2-naphthalenecarbonitrile (Enantiomer 1);
- 1-[(3,5-Dichlorophenyl)methyl]-3-[4-(4-fluorophenyl)-1-methyl-4-piperidinyl]-1*H*-pyrrole-2,5-dione;
- 1-[(3,5-Dichlorophenyl)methyl]-3-[4-(4-fluorophenyl)-1-methyl-4-piperidinyl]-5-methylidene-1,5-dihydro-2*H*-pyrrol-2-one; and phamaceutically acceptable salts thereof.
- 24. (Previously Presented) A compound according to claim 23 in amorphous or crystalline form.
- 25. (Previously Presented) A compound selected from:
- 1-[(3,5-Dichlorophenyl)methyl]-3-[4-(4-fluorophenyl)-1-methyl-4-piperidinyl]-1,5-dihydro-2*H*-pyrrol-2-one;
- 1-[(3,5-Dichlorophenyl)methyl]-3-[4-(4-fluorophenyl)-1-methyl-4-piperidinyl]-1,5-dihydro-2*H*-pyrrol-2-one hydrochloride;
- 1-[(3,5-Dichlorophenyl)methyl]-3-[4-(4-fluorophenyl)-1-methyl-4-piperidinyl]-1,5-dihydro-2*H*-pyrrol-2-one fumarate;
- 1-[(3,5-Dichlorophenyl)methyl]-3-[4-(4-fluorophenyl)-1-methyl-4-piperidinyl]-1,5-dihydro-2*H*-pyrrol-2-one citrate.
- 26. (Previously Presented) A compound according to claim 25 in crystalline form.
- 27. (Previously Presented) 1-[(3,5-Dichlorophenyl)methyl]-3-[4-(4-fluorophenyl)-1-methyl-4-piperidinyl]-1,5-dihydro-2*H*-pyrrol-2-one citrate.
- 28. (Previously Presented) 1-[(3,5-Dichlorophenyl)methyl]-3-[4-(4-fluorophenyl)-4-piperidinyl]-1,5-dihydro-2*H*-pyrrol-2-one.

- 29. (Previously Presented) A pharmaceutical composition comprising a compound as claimed in claim 1 in admixture with one or more pharmaceutically acceptable carriers or excipients.
- 30. (Previously Presented) A pharmaceutical composition comprising a compound as claimed in claim 25 in admixture with one or more pharmaceutically acceptable carriers or excipients.
- 31. (Previously Presented) A pharmaceutical composition comprising a compound as claimed in claim 28 in admixture with one or more pharmaceutically acceptable carriers or excipients.
- 32. 37. (Canceled).
- 38. 44. (Canceled).